

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX083122\
 Data File : VX031000.D
 Acq On : 01 Sep 2022 00:56
 Operator : JC/MD
 Sample : N4380-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-33

Quant Time: Sep 01 02:24:07 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X083022W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 30 14:24:53 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.556	168	221453	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	372140	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	346229	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	140228	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	138438	43.585	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.180%
35) Dibromofluoromethane	5.391	113	129528	47.218	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.440%
50) Toluene-d8	8.653	98	478244	47.075	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.160%
62) 4-Bromofluorobenzene	11.085	95	158850	43.909	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.820%
Target Compounds						
					Qvalue	
17) Carbon Disulfide	2.514	76	2435	0.343	ug/l #	79
19) Methyl tert-butyl Ether	3.111	73	4447	0.507	ug/l #	88
20) Methylene Chloride	2.788	84	1085	0.333	ug/l	88
30) Chloroform	5.105	83	13233	2.421	ug/l	89
40) Benzene	6.038	78	16600	1.413	ug/l	91
47) Bromodichloromethane	7.824	83	2269	0.551	ug/l	98
69) o-Xylene	10.646	106	1627	0.302	ug/l	92
95) Naphthalene	13.780	128	10027	0.816	ug/l #	92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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